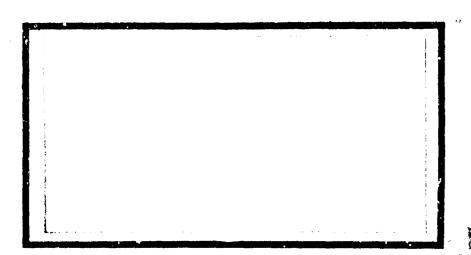
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DC	CONTROL DATA - R	& D	
(Security classification of title, body of ab	extract and indexing annotation must be e	intered when the	overall report is classified)
1. URIGINATING ACTIVITY (Corporate author)		24. REPORT SE	CURITY CLASSIFICATION
Air Force Institute of Te	chnology	Un	classified
Wright-Patterson Air Forc		26. GROUP	
	•	-	•
3. REPORT TITLE			
A MONTE CARLO TECHNIQUE U			
APPROXIMATE RELIABILITY C		SYSTEMS	WITH COMPONENTS
CHARACTERIZED RY THE WEIB	ULL DISTRIBUTION	•	•
4. DESCRIPTIVE NOTES (Type of report and Inclus-	ive dates)		
AFIT Thesis		· r · · · · ·	• • • • • • •
5- AUTHOR(\$) (First name, middle initial, last name	,)		
Robert G. Lannon		•	
Capt USAF			
Capt USAr		•	
6. REPORT DATE	74. TOTAL NO. O	F PAGES	7b. NO. OF REFS
March 1972	53	) )	15
88. CONTRACT OR GRANT NO.	Sa. ORIGINATOR'S	S REPORT NUME	BER(S)
	, 1		
b. PROJECT NO. N/A	GAM/MA/	72-2	
c.	9b. OTHER REPOR	RT NO(S) (Any of	ther numbers that may be assigned
	inte report,		
d.			
10. DISTRIBUTION STATEMENT			
American Community and an	12		
Approved for public relea	se; distribution uni	imitea.	
Approved for public release; IA	WAFR 190-127SPONSORING	MILITARY ACTI	VITY
Vocal A. W. Sleans	•		
Keith A. Williams, 1st Lt., US	AF [		
Acting Director of Information	-		

This paper develops a Monte Carlo technique which, with a digital computer, determines confidence limits for system reliability of complex systems containing components characterized by the Weibull distribution. The component distribution shape and scale parameters are estimated by the method of maximum likelihood from component failure times while the location parameter is assumed known. The asymptotic distribution of these maximum likelihood estimators and a Monte Carlo simulation are used to determine confidence limits on system reliability. As an example, confidence limits are calculated for two systems of up to eight components in combinations of series and parallel configurations using 99, 499, 999, and 2999 simulations. Accuracy of the confidence limits is found to be satisfactory after being checked by a method using a double Monte Carlo technique.

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13. ABSTRACT

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Security Classification

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KEY WORDS	LIN	K A	LIN	K B	LINK C		
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Monte Carlo Method				·			
Maximum Likelihood Estimators							
Reliability Estimates							
Weibull Density Functions							
Confidence Limits of Reliability							
Forecasting Reliability			-				
Computerized Simulation	!						
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A MONTE CARLO TECHNIQUE USING COMPONENT FAILURE TEST DATA TO APPROXIMATE RELIABILITY CONFIDENCE LIMITS OF SYSTEMS WITH COMPONENTS CHARACTERIZED BY THE WEIBULL DISTRIBUTION

GAM/MA/72-2

Robert G. Lannon Captain USAF



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A MONTE CARLO TECHNIQUE USING COMPONENT FAILURE
TEST DATA TO APPROXIMATE RELIABILITY CONFIDENCE
LIMITS OF SYSTEMS WITH COMPONENTS CHARACTERIZED
BY THE WEIBULL DISTRIBUTION

#### THESIS

Presented to the Faculty of the School of Engineering
of the Air Force Institute of Technology
Air University
in Partial Fulfillment of the
Requirements for the Degree of
Master of Science

by

Robert G. Lannon, B.M.S.

Captain USAF

March 1972

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### Preface

This thesis is a continuation of previous work done at the Air Force Institute of Technology on Monte Carlo techniques of reliability prediction. It is hoped that the method proposed here may find use in reliability analysis of complex systems and/or in situations where limited failure test data is available to the reliability engineer.

I wish to thank my thesis advisor, Professor Albert H.

Moore, for suggesting the topic and giving guidance and
encouragement throughout this project.

I also wish to express my gratitude to Dr. H. Leon Harter who gave valuable advise during this project. The FORTRAN programs in Appendix A and Appendix C are revisions of programs which he originally wrote and kindly supplied to me.

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#### Abstract

This paper develops a Monte Carlo technique which, with a digital computer, determines confidence limits for system reliability of complex systems containing components characterized by the Weibull distribution. The component distribution shape and scale parameters are estimated by the method of maximum likelihood from component failure times while the location parameter is assumed known. The asymptotic distribution of these maximum likelihood estimators and a Monte Carlo simulation are used to determine confidence limits on system reliability. As an example, confidence limits are calculated for two systems of up to eight components in combinations of series and parallel configurations using 99, 499, 999, and 2999 simulations. Accuracy of the confidence limits is found to be satisfactory after being checked by a method using a double Monte Carlo technique which assumes values for parameters of the component distributions and generates component failure times to be used in the proposed Monte Carlo technique. Central processing time for this technique on the CDC 6600 digital computer is found to be less than two minutes for the examples given. There is no limit to the number of components or the configurations of the components for the systems to which the method can be applied. The FORTRAN IV computer programs used are given along with flowcharts to facilitate reading the programs.

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### I. Introduction

#### The Problem

The purpose of this thesis is to determine whether accurate interval estimation of reliability for complicated systems, containing dissimilar components characterized by the Weibull distribution, can be obtained by using maximum likelihood estimates and a Monte Carlo simulation.

Significance. Systems containing many components become increasingly complex and costly. Complicated systems costing thousands of dollars can become inoperative due to the failure of a single connection. Because of this, it is important to determine system reliability as a function of the component reliabilities which make up the system. This can be done fairly easily for some few types of combinations of components which are from certain parent populations such as an electrical circuit made up of components characterized by the exponential distribution. When the components are dissimilar, no simple analytical expression can be used to compute system reliability.

In addition, predictions of reliability should be based on test data. However, as system reliability requirements

increase, testing a system to destruction can become expensive, both in time needed and money spent.

A method that circumvents this problem uses limited test data to calculate maximum likelihood estimates of the parameters of the failure distribution. These estimates are used, along with a random number generator, in a Monte Carlo technique which uses the asymptotic distribution of the maximum likelihood estimators to generate more point samples of system reliability. These point samples are used to determine an interval estimate of system reliability with the associated confidence level.

<u>Definitions</u>. Reliability is the probability that a system is still operating at time t. If T is the time to failure or life length of a system or component, the reliability at time t or R(t) is given by R(t) = P(T > t), where P means "probability of".

The Weibull density function is defined as

$$f(t;C,\theta,K) = \left[\frac{K(t-C)^{K-1}}{\theta^{K}}\right] \exp\left\{-\left[\frac{t-C}{\theta}\right]^{K}\right\}$$

$$\theta, K > 0, C \le t$$
(1)

#### = 0 elsewhere

where  $\theta$  = the scale parameter, K = the shape parameter, and C = location parameter. The scale parameter affects the dispersion of the random variable t about its mean. The shape parameter determines whether the hazard function is

increasing, decreasing, or time invariant; while the location parameter determines the point of origin.

Reliability, for the Weibull density function, is

$$R(t) = \exp \left\{ -\left[\frac{t-C}{\theta}\right]^{K}\right\}$$
 (2)

Pseudo-random numbers are generated from a series of numbers which repeats itself with long periods between repetitions. These numbers are suitably random if the repetition periods are greater than the number of digits desired and if there is a small enough correlation between digits (Ref 15:257).

The maximum likelihood estimator (NLE) is the estimator which maximizes the likelihood function with respect to the parameter being estimated. The likelihood function is the joint density of a sample of n random variables and is expressed by

$$L(x_1,x_2..x_n;\theta) = \prod_{i=1}^{n} f(x_i;\theta)$$
 (3)

where  $L(x_i:\theta)$  = the likelihood function.

If it is desired to determine  $\hat{\theta}$  (the MLE for the parameter  $\theta$ ), normally it is only necessary to take the partial derivative of the likelihood function with respect to  $\theta$ , set it equal to zero and solve for  $\theta$  which then becomes  $\hat{\theta}$ . By using the MLE for each parameter and the expression for reliability for a given failure density, it is possible to arrive at a point estimate for reliability.

The method in this paper generates intervals from a random sample of system reliabilities and this is called a random interval. If reliability falls within any given interval, there is associated with the interval a confidence level. For example, if it is said the reliability of a system lies between 0.6 and 0.8 at the 95% confidence level, it is meant  $P(0.6 \le R(t) \le 0.8) = .95$  or the random interval 0.6 - 0.8 would contain the true system reliability 95% of the time if the system were tested many times. The interval is called a confidence interval, which has 0.6 as a lower confidence limit and 0.8 as an upper confidence limit.

A Monte Carlo simulation is one in which a random value or number is picked from all possible numbers that can be described by a specific density function.

#### Analysis

Assumptions. This thesis will concern itself with systems which have components described by the Weibull density function. It is assumed that the components have previously been determined to be Weibull or that the Weibull density function adequately models the components in the system. This paper does not provide a method to test for the Weibull density function though these tests do exist and are described in other references. One of the easiest ways to test for the Weibull distribution is to use Weibull graph paper (Ref 5:170-173). Once it is determined or suspected the components are Weibull, the methods in this paper can be

used to estimate the parameters of the Weibull distribution and then to generate interval estimates of system reliability. The Weibull distribution is rather versatile in that, depending on the value of the shape parameters, the Weibull distribution can be used to model components with decreasing, increasing or constant failure rates or hazard functions.

A special case of the Weibull distribution is the exponential distribution which occurs when the shape parameter, K, is equal to one.

It is assumed that the components of the system being analyzed fail independently. This allows use of standard formulas for components connected in series or in parallel. For instance, if a system is composed of two components connected in series, the reliability can be expressed by

$$R_s(t) = [R_1(t)][R_2(t)]$$
 (4)

where  $R_1(t)$  = reliability of component number one and,  $R_2(t)$  = reliability of component number two.

Likewise, for a system made up of two components in parallel the reliability can be expressed by

$$R_s(t) = 1 - [1 - R_1(t)][1 - R_2(t)]$$
  
= 1 - Q<sub>1</sub>Q<sub>2</sub> (5)

where  $Q_i = 1 - R_i(t)$ .

More complex systems can be reduced to combinations of series and/or parallel configurations by use of Bayes'

theorem or the Boolean Disjunctive theorem. In other cases where the system reliability dependence on component reliability is known and can be expressed by other equations, the independent failure assumption is not necessary.

It is assumed that the location parameter is known or can be set equal to zero for all Weibull component distributions in this paper.

Standards. The only true criterion that is applied to test the acceptability of the method in this paper is analysis of the result. Does the computed confidence interval contain the actual value of system reliability the required number of times? To check the validity of the method, the values of all parameters were assumed for each of the Weibull distributions of the components. Pseudo-random component failure times were generated for each of the components and the method in this paper applied to generate many confidence intervals. Since the true system reliability was known, a check was made to see if the true reliability was in the calculated intervals the required number of times. This process is described in Chapter II.

For the estimation of the parameters from the generated component failure times, variation from the true value of the parameter in the second digit was deemed acceptable.

Background. Levy and Moore developed a method for determining system reliability confidence limits using a Monte Carlo method where the component failure distributions (exponential, normal, lognormal, Gamma, or Weibull) with

location and shape parameters were known (Ref 10). Lutton extended this method by using the asymptotic distribution of the maximum likelihood estimators of, among others, the Weibull distribution, location parameter known (Ref 11).

Harter and Moore developed two procedures relevant to this paper. They used an iterative procedure to determine maximum likelihood estimates of the parameters of the Weibull distribution from complete and censored samples (Ref 8). They also developed a method to determine the asymptotic variances and covariances of the maximum likelihood estimators of the parameters of the Weibull distribution from samples that were censored or complete (Ref 7). This paper presents a technique in which the sample of failure times for each component is used, along with the asymptotic distribution of the Weibull parameters, to generate a confidence interval with the associated confidence level.

Approach. The method and associated computer program in this paper:

- Calculate the MLE of the component shape and scale parameters from the component failure times.
- Determine the variance-covariance matrix of the asymptotic normal distribution of the shape and scale parameters.
- Generate a sample of the component shape and scale parameters.
- 4. Use sample shape and scale parameters, along with known location parameter, to calculate a point

estimate of component reliability for each component.

- 5. Calculate a point sample of system reliability from the point samples of component reliabilities.
- 6. Repeat steps 3-5 to obtain many samples of system reliability.
- 7. Order the system reliability samples to obtain an interval sample of system reliability from which the confidence limits are determined at a given confidence level.

#### II. Procedure

#### Estimation of Component Parameters

The computer program for finding the MLE of the parameters of the Weibull distribution from complete or censored samples is given in Appendix A. The procedure, as developed by Harter and Moore, uses the natural logarithm of the likelihood function of the remaining order statistics which is

$$L_r = \ln n! - \ln(n-m)! - \ln r! + (m-r)(\ln K - K \ln \theta) +$$

$$(K-1)\sum_{i=r+1}^{m} \ln(x_i-C) - \sum_{i=r+1}^{m} \left[\frac{(x_i-C)}{\theta}\right]^{K} -$$

$$(n-m)\left[\frac{x_m-C}{\theta}\right]^K + r \ln\left\{1 - \exp\left[\frac{-(x_{r+1}-C)^K}{\theta^K}\right]\right\}$$
 (6)

where (m-r) is the number of order statistics  $x_{r+1}$ ,  $x_{r+2}$ ,..., $x_m$  and n is the size of the sample before censoring.

The partial derivatives of  $L_r$  with respect to each of the three parameters are set equal to zero and the three resulting equations solved simultaneously. The iterative procedure for the simultaneous solution of the equations uses an initial estimate of the parameter (which is fed into the computer along with the failure times) and the rule of false position to determine a new value of the parameter from the appropriate likelihood equation into which the latest estimates (or a known value) of the other two parameters have been substituted (Ref 8:641).

The program in Appendix A has been converted into a subroutine and is called for analysis of each component. When this is done, the MLE of the shape and scale parameters are determined for all components in the system.

### Generating Sample System Reliabilities

The asymptotic joint distribution of the estimators with the location parameter known is bivariate normal and is given by

$$f(Y) = \frac{1}{|V|^{1/2}(2\pi)} \left[ \exp\left(-\frac{1}{2} Y'V^{-1}Y\right) \right]$$
 (7)

where Y = two component vector of estimates of the parameters

Y' = transpose of Y

V = asymptotic variance-covariance matrix

 $V^{-1}$  = inverse of V

|V| = determinant of V

Let  $Z = V^{-1/2}Y$  so that  $Y = V^{1/2}Z$  and, because V is symmetric,  $Y' = Z'V^{1/2}$ . Then

$$\hat{x}(?) = \frac{1}{2\pi} \left[ \exp\left(-\frac{1}{2} z' v^{1/2} v^{-1} v^{1/2} z\right) \right]$$
 (8)

$$= \frac{1}{2\pi} \left[ \exp\left(-\frac{1}{2} Z'Z\right) \right]$$
 (9)

$$= \frac{1}{2\pi} \left[ \exp \left( 2_1^2 + 2_2^2 \right) \right] \tag{10}$$

so that the density function of Z is made up of independent variates of the normal distribution with mean zero and variance 1 (Ref 14). Thus, to obtain a sample of the joint distribution of the estimators, the following formula is used:

$$Y' = V^{1/2}Z + Y$$
 (11)

where Y' = 2 X 1 vector of the biased random samples of the estimates of the shape and scale parameters.

V = 2 X 2 variance-covariance matrix.

Z = 2 X 1 vector of pseudo-random numbers from the standard normal distribution.

Y = 2 X 1 vector of the maximum likelihood estimates

of the two parameters (as found in the previous

section).

or

$$\begin{pmatrix}
\hat{\theta}_{\text{sample}} \\
\hat{\kappa}_{\text{sample}}
\end{pmatrix} = \begin{pmatrix}
v_{11} & v_{12} \\
v_{21} & v_{22}
\end{pmatrix} \begin{pmatrix}
z_{1} \\
z_{2}
\end{pmatrix} + \begin{pmatrix}
\hat{\theta}_{\text{MLE}} \\
\hat{\kappa}_{\text{MLE}}
\end{pmatrix} (12)$$

In formula (12) the MLE of  $\theta$  and K are used as the marginal means of the bivariate normal distribution of the parameter estimates.

The random samples of the component parameter estimates were substituted into the expression for reliability for each component to get random values of component reliabilities. The were then substituted into the expression for system

reliability, one value for each component, to generate random values of system reliability.

The program which generated the system reliabilities is given in Appendix B, along with a flow chart of the program. A detailed explanation of the evaluation of  $V^{1/2}$  is also given in Appendix B.

### The Variance-Covariance Matrix

The natural logarithm of the likelihood function for a sample of size n from a Weibull population where the lowest r and the highest n-m sample values have been censored is given in Ref 7. The elements of the information matrix are found by taking the limits, as  $n \to \infty$ , of the negatives of the expected values of the second partial derivatives of the likelihood function with respect to each of the parameters. Let  $k_{ij}$  be the elements of a matrix B, formed from the elements of the information matrix  $(s_{11}, s_{12} = s_{21}, and s_{22})$ , where E[...] is the conditional expectation given  $z_{r+1}$  and  $z_m$ . Then

$$k_{11} = \frac{v_{11}\theta^{2}}{n} = \lim_{n \to \infty} \Pr \left[ -\left(\frac{\theta^{2}}{n}\right) \left(\frac{\partial^{2}L}{\partial\theta^{2}}\right) \right]$$

$$= -Kp + K(K+1) \left[\Gamma(2;\hat{\tau}_{m}^{K}) - \Gamma(2;\hat{\tau}_{r+1}^{K})\right] + K(K+1)q_{2}\hat{\tau}_{m}^{K} + \frac{\hat{\tau}_{r+1}f(\hat{\tau}_{r+1})[K\hat{\tau}_{r+1}^{K} - (K+1)q_{1}]}{q_{1}}$$
(13)

$$k_{12} = \frac{v_{12}\theta}{n} = \lim_{n \to \infty} \operatorname{pr} E \left[ -\left(\frac{\theta}{n}\right) \left(\frac{\partial^{2}L}{\partial\theta\partial K}\right) \right]$$

$$= p - \left[\Gamma'(2;\hat{z}_{m}^{K}) - \Gamma'(2;\hat{z}_{r+1}^{K})\right] - \left[\Gamma(2;\hat{z}_{m}^{K}) - \Gamma'(2;\hat{z}_{m}^{K})\right] - \left[\Gamma(2;\hat{z}_{m}^$$

$$k_{22} = \frac{v_{22}}{n} = \lim_{K \to \infty} \operatorname{pr} \left[ -\left(\frac{1}{n}\right) \left(\frac{\partial^{2} L}{\partial^{2} K}\right) \right]$$

$$= \frac{p}{K^{2}} + \frac{\left[\Gamma''(2; \hat{z}_{m}^{K}) - \Gamma''(2; \hat{z}_{r+1}^{K})\right]}{K^{2}} + q_{2}\hat{z}_{m}^{K} \ln^{2} \hat{z}_{m} + \frac{\hat{z}_{r+1}^{K} + \hat{z}_{r+1}^{K}}{Kq_{1}} \left\{ \ln^{2} \hat{z}_{r+1}^{K} + \frac{\hat{z}_{r+1}^{K} - q_{1}}{Kq_{1}} \right\}$$
(15)

where  $q_1 = r/n$ ,  $q_2 = (n-m)/n$ , and  $p = 1 - q_1 - q_2 = (m-r)/n$ . As  $n \to \infty$  with  $q_1$  and  $q_2$  fixed,  $z_{r+1}$  converges in probability to  $\hat{z}_{r+1}$  where  $F(\hat{z}_{r+1}) = \int_0^{\hat{z}_{r+1}} f(t)dt = q_1$  and  $z_m$  converges in probability to  $\hat{z}_m$  where  $1 - F(\hat{z}_m) = \int_{z_m}^{\infty} f(t)dt = q_2$ . The primes indicate differentiation and  $\Gamma(a;b)$  is the incomplete Gamma function (Ref 7:559).

Then the information matrix, in terms of the elements  $\mathbf{k}_{\mathbf{i}\mathbf{j}}$ , becomes

$$\begin{pmatrix} s_{11} & s_{12} \\ s_{21} & s_{22} \end{pmatrix} = \begin{pmatrix} \frac{k_{11}^n}{\theta^2} & \frac{k_{12}^n}{\theta} \\ \frac{k_{21}^n}{\theta} & nk_{22} \end{pmatrix}$$
(16)

Let  $a_{ij}$  be the elements of the matrix A, which is the inverse of B, then

$$\begin{bmatrix} A \end{bmatrix} = \begin{bmatrix} a_{11} & a_{12} \\ a_{21} & a_{22} \end{bmatrix} = \begin{bmatrix} k_{11} & k_{12} \\ k_{21} & k_{22} \end{bmatrix}^{-1} = \begin{bmatrix} B \end{bmatrix}^{-1}$$
 (17)

The elements of the variance-covariance matrix are then

$$\begin{pmatrix} v_{11} & v_{12} \\ v_{21} & v_{22} \end{pmatrix} = \begin{pmatrix} \frac{a_{11}\theta^2}{n} & \frac{a_{12}\theta}{n} \\ \frac{a_{21}\theta}{n} & \frac{a_{22}}{n} \end{pmatrix}$$
(18)

since

$$\begin{pmatrix} \frac{k_{11}^{n}}{\theta^{2}} & \frac{k_{12}^{n}}{\theta} \\ \frac{k_{21}^{n}}{\theta} & k_{22}^{n} \end{pmatrix} \begin{pmatrix} \frac{a_{11}^{\theta^{2}}}{n} & \frac{a_{12}^{\theta}}{n} \\ \frac{a_{21}^{\theta}}{n} & \frac{a_{22}^{\theta}}{n} \end{pmatrix} = \begin{pmatrix} k_{11}^{a_{11}} + k_{12}^{a_{21}} & \frac{1}{\theta} & (k_{11}^{a_{12}} + k_{12}^{a_{22}}) \\ \theta (k_{21}^{a_{11}} + k_{22}^{a_{21}}) & k_{21}^{a_{12}} + k_{21}^{a_{22}} \end{pmatrix} = I \quad (19)$$

The program given in Appendix C finds the elements of the asymptotic variance-covariance matrix by solving Eqs (13),

(14), and (15) to find the elements  $k_{ij}$ , inverting them and converting these to the elements of the variance-covariance matrix by the above transformations.

The variance-covariance matrix is solved by this program for non-integer values of the shape and scale parameters since with the location parameter known the equations are regular for any non-integer value of the parameter estimates. A flowchart and more detailed explanation of the equations in the program are also included in Appendix C.

### Calculating System Confidence Limits

The sample values of system reliability, after having been generated, are ordered and thus yield the sample cumulative distribution of the system reliability. From this distribution, the confidence interval and limits can be approximated at any level of confidence. For example, in an ordered sample of 99, the tenth value represents the lower limit of a one sided confidence interval at the 90% confidence level since the a priori probability is .90 that a new random value exceeds the tenth order statistic of a sample of 99. Thus, it is assured that the true reliability is equal to or greater than this value with a confidence of .90.

### Verifying the Level of Confidence

In order to insure that the confidence level is accurate, a program was written to generate sample times to failure from the Weibull distribution. Parameters were assumed for each of the component distributions and various size samples

of time to failure were generated for each of the components. This simulates bench testing each of the components to obtain a sample of failure times for each component. The computer program which does this is given, with an explanation of the procedure, in Appendix D.

The <u>true</u> reliability of the system is now known, since the component distribution parameters are known.

The failure times for each component were fed into the Monte Carlo program and a number of sample values of the confidence interval at the required level were generated. The true reliability should be covered by a percentage of the sample confidence intervals equal to the confidence level. For example, if 10 one sided 50% confidence intervals are generated and the true system reliability is .90, a graph of the intervals might appear as shown in Fig. 1 on the next page.

In this case half the intervals cover the true reliability verifying the 50% confidence level (Ref 13:253).

For this paper confidence intervals were generated with each interval calculated from a sample (N) of 99, 499, 999, and 2999 system reliabilities. The results are given in Chapter III.

#### Systems Analyzed

Two systems are analyzed in this paper. The first system consists of 8 components connected as shown in Fig. 2 for which the required mission time is 75 hours.

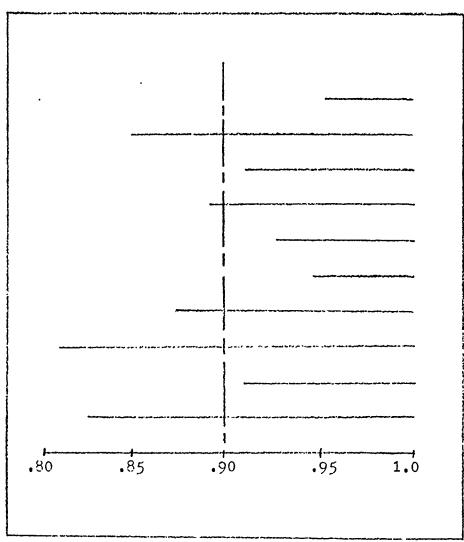


Fig. 1. A Sample of Ten Confidence Intervals Over a True Reliability of .90

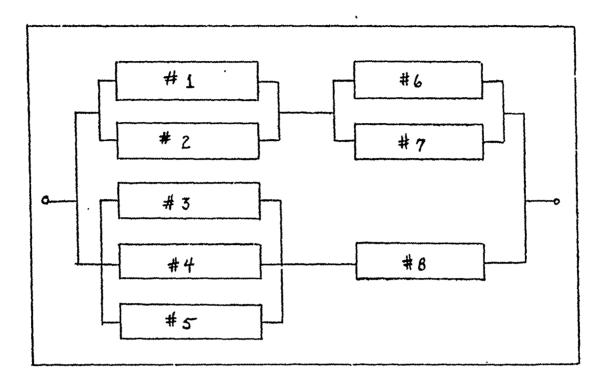


Fig. 2. Schematic of System A.

The components have the following characteristics:

## Component #1.

Failure distribution: Exponential

Failure test sample size: 100

Sample censoring: top 6 and bottom 6 values

Location parameter: 0.0

### Component #2.

Failure distribution: Exponential

Failure test sample size: 75

Sample censoring: None

Location parameter: 0.0

### Component #3.

Failure distribution: Weibull

Failure test sample size: 150

Sample censoring: None

Location parameter: 0.0

### Component #4.

Failure distribution: Weibull

Failure test sample size: 200

Sample censoring: None

Location parameter: 0.0

### Component #5.

Failure distribution: Exponential

Failure test sample size: 75

Sample censoring: None

Location parameter: 0.0

### Component #6.

Failure distribution: Weibull

Failure test sample size: 100

Sample censoring: None

Location parameter: 0.0

### Component #7.

Failure distribution: Weibull

Failure test sample size: 200

sample censoring: None

Location parameter: 0.0

#### Component #8.

Failure distribution: Weibull

Failure test sample size: 250

Sample censoring: top 12 and bottom 12 values

Location parameter: 0.0

The expression for the reliability of this sastem is

$$R_s(t) = 1 - Q_a(t)Q_b(t)$$
 (20)

where  $Q_i = 1 - R_i(t)$   $R_a(t) = [1 - Q_1(t)Q_2(t)][1 - Q_6(t)Q_7(t)]$   $R_b(t) = [1 - Q_3(t)Q_4(t)Q_5(t)]R_8(t)$ 

The second system (system B) has the configuration shown in Fig. 3 with a required mission time of 100 hours.

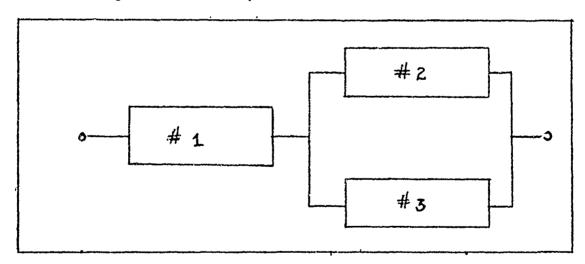


Fig. 3. Schematic of System B.

The components of system B have the following characteristics:

### Component #1.

Failure distribution: Weibull

Failure test sample size: 10

Sample censoring: None

Location parameter: 0.0

### Component #2.

Failure distribution: Weibull

Failure test sample size: 20

Sample censoring: None

Location parameter: 0.23

### Component #3.

Failure distribution: Exponential

Failure test sample size: 30

Sample censoring: None

Location parameter: 0.0

The expression for system reliability for system B, in terms of the  $i^{\mbox{th}}$  component reliability is:

$$R_s(t) = R_1(t)(1 - Q_2(t)Q_3(t))$$
 (21)

System B was designed to put a strain on the proposed Monte Carlo method due to the small failure test sample size for each of the components.

#### III. Results

### Confidence Limits

It is known that the error associated with a Monte Carlo calculation is proportional to  $1/\sqrt{N}$  where N is the number of simulations (Ref 15:259). In this case, the error is statistical, that is, the probable error is proportional to  $1/\sqrt{N}$  or the probability is high that the approximate solution does not deviate from the true solution by more than a certain amount (Ref 12:255). The use of a high speed digital computer should decrease the amount of error in the calculation but the machine can add random errors of its own, such as round-off error (Ref 9:11,12). As the problem becomes more complex, the error can be estimated only from the results of the computation.

Tables I and II, on the following page, show the confidence limits determined from simulations of various size for systems A and B. As the simulation size increases the accuracy of the confidence limits at a given confidence level increases and more closely approximates the true confidence limits. Also given are the central processing times for the CDC 6600 computer for each size simulation. No trend should be noted from the bounds given for each size simulation in these tables since each simulation is a random sample of system reliabilities for a single run. The only proper statement that can be made is that the confidence limits for a simulation of size 2999 are more accurate than for one of

Table I
One Sided Confidence Limits
System A

System A													
% Conf	, <del> </del>	99	N=	=499	N=	999	N=2	5000					
Ievel	(a)*	('n)**	(a)	(ħ)	(a)	(b)	(a)	(5)					
99	1	•9739	5	.9731	20	.9744	30	•9735					
98	2	•9755	10	.9740	30	•9754	60	9745					
97	3	•9760	15	•9748	40	.9761	30	.9754					
96	4	•9762	20	•9756	50	. •9764	120	.9761					
95	5	.9752	25	•9762	60	.9768	i : 150 !	•9763					
90	10	.9772	50	•9775	100	•9779	i 300	•9777					
: : 85 . :	<b>1</b> 5	•9779	75	•9789	150	.9789	. 450 :	. <sup>072</sup> 6					
: 80 :	20	.9786	100	.9796	200	•979°	600 ;	• 62.09					
70	30	.9801	150	.9807	300	.9808	930	.9806					
60	40	.9810	200	.9815	400	.9816	1200	.9815					
50	50	.9816	250	.9323	500	.9824	1500	•òes3					
CP Time	19	.4 Sec	23.	0 Sec	29	.5 Sec	89.6 <b>S</b> ec						

<sup>\*</sup> Order number of the sample \*\* Confidence limit on system reliability

Table II One Sided Confidence Limits System B

% Conf	N:	<b>-</b> 99	V:	490	.i	999	N=2099			
Level	(a)* (b)**		(a)	(b)	(2)	(b)	(a)	(6)		
99	1	.8788	5	.8680	20	.8427	30	.8287		
98	2	•9077	10	.8888	30	.3632	60	.8669		
97	3	•9094	<b>1</b> 5	.8994	<u>40</u>	.8783	90	.3822		
96	Ļ	.9125	20	.9046	50	.8864	120	.2919		
95	5	.9134	25	.9077	60	.8961	150	.9014		
90	10	.9230	50	.9224	100	.9213	350	.9223		
85	15	•9357	75	•932h	<b>1</b> 50	•933# •	: 50	•9339		
80	20	.9388	100	•9395	200	*uħ 58	600	.9-12		
70	. 30	•9487	<b>1</b> 50	•9489	300	.9524	900	.9518		
60	10	•9549	200	•9558	400	•9576	1200	•9575		
50	50	.9616	250	.9605	500	.9626	1500	.9628		
CP Time	13.	2 Sec.		.4 Sec.		.3 Sec.	76.	O Sec.		

<sup>\*</sup> Order number of sample \*\* Confidence limit on system reliability

size 99. Only the single sided confidence limits are given in these tables because they are the numbers most needed in evaluating the reliability of a new system.

There is also an error associated with the size of the sample of failure times for each component. With the location parameter known, the sample variances and covariances of the asymptotic distribution of the shape and scale parameters are different from the variances and covariances of the small-sample non-asymptotic distribution of the parameters.

Analyzing this error is difficult but Harter and Moore determined the error to be relatively small if the location parameter is known, even for samples of failure times as small as size 50 (Ref 7:562, 563).

Lutton compared the asymptotic distribution technique with a double Monte Carlo simulation developed by Moore to approximate the system reliability confidence interval.

Moore's method does not use the asymptotic distribution of the parameters but uses a method that has an accuracy that is dependent upon the number of simulations made. Lutton found good agreement between the asymptotic method and the double Monte Carlo method (Ref 11:21-26).

### Verification of Confidence Level

In order to check the accuracy of the confidence level, the proposed method was repeated several times. Here, a single repetition of the method is called a run. At first, because of the computer time needed for a single run for

system B, 100 runs were made on both systems A and B for a sample of 99 system reliabilities (simulation size 99). In each run the following is accomplished:

- Sample failure times for each component are generated from component Weibull distributions with assumed parameters.
- An ordered sample of 99 system reliabilities is calculated by the proposed method.
- 3. A counter is increased by one if the true reliability of the system is in the generated confidence interval at any given confidence level.

After all the runs are completed, the value of the counter is divided by the total number of runs. This gives the percentage of runs in which the true reliability is in the confidence interval at a given level; as in the example shown in Fig. 1. The results of this procedure are given in Table III on the next page. As shown in this table the confidence level is found to be reasonably accurate for a simulation size 99, which is the least accurate interval because of the error associated with that size sample.

Table III indicates that the established or required confidence levels in the proposed Monte Carlo method are sometimes optimistic, at other times pessimistic. It must be remembered that these confidence levels were checked by using only 100 runs on a simulation size 99 and it can be expected that the calculated levels would approach the required levels if more runs were possible. Even with these limitations,

Table III Calculated Confidence Levels N=99\*:100 Runs

Required Conf. level	Calculated (7)		
	System A	System 3	
·99S**	99	99	
9°T#	99	95	
905	85	<b>6</b> , <del>1</del>	
òoū	en	àδ	
808	75	86	
80T	76	96	
708	62	75	
?0型	64	79	
608	56	57	
60т	59	72	
508	46	45	
50Т	46	62	

<sup>\* 99</sup> system reliabilities simulated\*\* One sided confidence interval# Two sided confidence interval

Table III indicated the accuracy of the proposed method is fairly good at a given confidence level.

To determine the effect of more runs on the calculated confidence level. System B confidence levels were calculated for 300 and 700 runs. This was done only for System B because computer time for each run is not excessive for this three-component system. The results are given in Table IV on the next page. In most cases, if the calculated level was below the required level, the calculated level approaches the required level and then surpasses it when more runs are made. If the calculated confidence level was above the required level, the calculated level approaches the established or required level when more runs are made. In almost all cases the required confidence levels for this system tend to be conservative. This confirms Lutton's results, ir which he found the confidence intervals for the asymptotic distribution of the parameters tended to be wider or more conservative than in Moore's double Monte Carlo method. In other words, if the results in Table IV apply to any system analyzed, the confidence limits in the proposed Monte Carlo method tend (in a probabilistic sense) to be slightly conservative for both one and two sided confidence intervals. This is a desirable characteristic from a practical standpoint.

Table IV System B Confidence Levels with Runs Increased N=99\*

Required Conf. Level	Calculated (%)			
	100 Runs	300 Runs	700 Runs	
995**	99.0	99.0	99.5	
98T #	95.0	96.0	97.3	
908	94.0	92.3	92.6	
90Т	88.0	89.0	90.4	
80S	86.0	80.6	£2 <b>.</b> 6	
80T	86.0	82.3	83,6	
705	76.0	72.0	73.0	
70T	79.0	72.3	74.1	
605	57.0	57.0	60.4	
60т	72.0	63.3	64.4	
50\$	45.0	45.0	49.0	
50Т	62.0	54.6	53•9	

<sup>\* 99</sup> system reliabilities simulated\*\* One sided confidence interval# Two sided confidence interval

# IV. Conclusions and Recommendations

### Conclusions

The method developed is extremely flexible and gives good accuracy for a wide number of system configurations.

The only inputs needed to start the process are the ordered censored or uncensored samples of component failure times and, along with the known value of the component location parameters, any kind of estimates of the component scale and shape parameters. In the latter case a guess will suffice since these estimates are used only to start the method of false position. An increasing or decreasing failure rate may be indicated by the failure times, thus supplying a guess for the shape parameter. Any number within the range of failure times may be used to guess the scale parameter.

Bayesian methods or Boolean algebra may be used to express system reliability in terms of the component reliabilities. In most cases this can be done with a single equation which is inserted in the main program. The proposed method can be applied to any component combination and uses relatively little computer time to generate a large sample from which confidence limits are determined.

The accuracy of the method depends mainly on the number of simulations of system reliability. It also depends, to a lesser extent, on the size of the sample of component failure times. As the size of either of these increases so does the accuracy of the confidence interval. The confidence limits

are easily recovered from the ordered sample of system reliabilities printed as output from the computer program.

### Recommendations

It is recommended that further work be done on the proposed Monte Carlo method in the following areas:

Types of Distributions. The method should be extended to include other distributions such as the Logistic and Gamma distributions. Iterative procedures have been developed to obtain the MLE for the parameters of these distributions and could be incorporated into a program such as the one in this paper.

Error Analysis. It is also recommended that a wider variety of component configurations be analyzed to try to determine if there is a direct relationship between the number and configuration of the components and the error. In other words, is error prediction possible, or can error be determined only by analysis of the results?

Analysis of systems with component failure time samples of less than ten would also be useful. If the proposed method remains accurate for such limited component testing, it could be used in systems which have components with a long mean time to failure or where the components are so expensive that only a limited number could be tested to destruction.

Confidence level verification as given in Table III uses a limited number of computer runs on simulation size of 99 system reliabilities. It is recommended that further examination of the accuracy of the confidence level for simulations

of 499, 999, and 2999 also be undertaken to check the increase in accuracy of the confidence interval for these larger size simulations.

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### Appendix A

# The Parameter Estimation Subroutine

```
SUPROUTING PARES (N, M, C, TH, TA, FK, MR, FTH, PEK)
                                                      THE THE CONTROL OF CHASORING, A = 100 OF LESS AS

STATE IF SULL BARAMETER THATA IS KNOWN

SCIENT IF SULL BARAMETER THATA IS TO BE ESTIMATED

SCIENT IF SULL BARAMETER K TO BE ESTIMATED

SCIENT IF SUL
                                                           gurruit
                                                       SS 2= 1.
SS 2= 1.
SS 2= 1.
SF (%) 55,56,136
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          1 4 [N=N

PEA) 3, (T(I), T=1, N)

3 FORWAT (5F16.5)

IF (*) **5.56, -1

11) PRINT 115

115 FURVAT(14 ,/, IH ,15HESTIMATES!)

PRINT 121

120 FORWAT (14 , HUCCATTON, TX, SHECKLE
         ---- 11 St)

---- 12 St)

---- 
           PRINT 125
125 FORMAT(14 ,//,1 + ,1 ,41T+ RATIONE)
                                                         EM=>".
 31
                                                           K 42= 42+1
                                                          NY=K-M+1
DO [4 T=\M,1]
33
                                                          FIFI
                                                         ELAV=FLKM+/LOG(II)
IF (MP) 66,15,75
DO /F I=4,40
 34
75
                                                          FERYERINY-NLCG(TT)
```

```
DO ?. J=1,551
IF (J-1) 21,27,77
35
     JJ=J-1
SK= •
37
     SL=^.
     St=0.
DC : I=var.w
Sk=Sk+(T(I)-C(JJ))**FK(JJ)
     IF (St1) 7,7,8
THEIN(J)=THEIN(JJ)
GO TO 0
6
7
      IF (YP) 66,19,2
THETA(J)=((5K+(5N+F5)*(T(M)+5(JU))**FK(JU))/EM)
A
    C##(1./~K(JJ))
GC TO G
X(1)=THTTS(JJ)
23
     ËS=
     50 21 L=1,5
      LL=[-1
      [[=]+1
    えんじょうきん(じ)
53
     Ĭ: () C
54
     IF ([S=1]) [], 35, 58
X([C) = .55 Y([)
GO TO 51
                                           Reproduces from available copy.
55
      Y(LT)=1.5*Y(L)
55
     58
50
6.6.2.7
      9
11
17
      X(1)=FK())
L9=
      $FK=PFK+(FFU@(L(I)+C(AA))+VFUG(LH=IV(A)))*(L(I)+C(AA));

BFK=J*

BO_1 F=T*>U
13
      LL=1-1
    LP=1+1
47
44
45
45
      IF (Y(L) #Y(LL)) 47, ~7,47
47
43
      1.L=LL-1
```

1

The state of the s

```
X(Lr):X(L):Y(L):*(Y(L)-X(LL))/(Y(LL)-Y(L))
IF:(ARS(Y(LP)-Y(L))-1:L-4) 5?,%_,51
CONTINUT
5516171
                                                        UNITER (1) = ((LF)

C(J) = ((JJ)

IF (SS3) 25,25,14

IF (1,-74(J)) 15,73,78

IF (SS1+3S2) 57,57,16

X(1) = C(J)

LS=1
                                                          DO 23 L=1.55

SK1=7.

SR= .

DO 15 l=vrc.v

SK1=SK1+(*([])-v([))**(f<([])-1.)

SP=SR+1./(*([])-v([))
  15
                                            LL=1+1

X(LC)=Y(L)

ZPK=((T(MAC)-Y(L))/TH(T'(J))**TK(J)

Y(L)=(1.-1x(J))*CP4+(Y(J))*(CX1+(-h-2Y)*(T(Y)-Y(L)))

C**(TX(J)-1.)/JUCTA(J)**TX(J)-1MP*CX(J)-7MFFXC(-7CY)

C/((T(MAC)-Y(L))*(1.-2*C(-2?*)))

LC=(X(J)-1.)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/CP+(L)/
                                                             LL=1 ~1
     37
   43
      41
      42
                                                               X(LT)=X(L)+Y(L)*(Y(L)-Y(LL))/(Y(LL)-Y(L))
IF (ATC(X(LT)-Y(L))-1.--4) 24,24,23
CONTTACT
      23
                                                             CONTINUE
C(J)=Y(LP)
GC TC TC
C(J)=T(1)
TF (MP) Eq. 18.33
DO SE 1=1.4
IF (C(J)+1.7-4-T(I)) 69.67,57
MT=MT+1
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                           Reproduced from sold best available copy.
       25
                                                               Mn= (2) +1 - 4 - C(1) = T(1) | IF (40) | 50,79,71
       67
63
68
                                                  IF (\(\frac{\partial}{\partial}\) 36,(3,\frac{\partial}{\partial}\)
SL=\(\frac{\partial}{\partial}\)
SL=\(\frac{\partial}{\partial}\)
SK=\(\frac{\partial}{\partial}\) (J)\) ***\(\frac{\partial}{\partial}\)
SL=\(\frac{\partial}{\partial}\) (T(\frac{\partial}{\partial}\) (J)\) ***\(\frac{\partial}{\partial}\) (J)\) ***\(\frac{\partial}{\partial}\) (J)\) ***\(\frac{\partial}{\partial}\) (T(\frac{\partial}{\partial}\) **\(\frac{\partial}{\partial}\) (T(\frac{\partial}{\partial}\) **\(\frac{\partial}{\partial}\) (T(\frac{\partial}{\partial}\) **\(\frac{\partial}{\partial}\) (T(\frac{\partial}{\partial}\) **\(\frac{\partial}{\partial}\) (T(\frac{\partial}{\partial}\) **\(\frac{\partial}{\partial}\) 
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        36
          27
28
                                                           50
           3.)
                     125
136
                      127
                                                      Chall
```

PIH=IH=TA(J)

PEK=FK(J)

155 FORMAT (14)

140 CONTTNUE

KETLEN

END

### Appendix B

# Generating the System Reliabilities

The main Fortran program SYSRE generates the system reliabilities. The inputs to this program are the number of components in the system (a maximum of ten as presently dimensioned), the number of simulations desired, and the mission time. Censoring information is supplied to the program along with the known value of the location parameter. The failure times are read for each component when the subroutine PARES is called and the true value of the location parameter is read into the main program for each component. The output from the main program is a sample of ordered system reliabilities from which the confidence limits can be read for a given confidence level. The flowchart is given in Fig. 4 and the function subprogram which is used to generate pseudo-random numbers from the standard normal distribution is explained in the notes.

# Notes (refer to Fig. 4):

- 1. The ordered component failure times are read into subroutine PARES and the individual component MLE's of the shape and scale parameters are calculated and stored each time the subroutine is called.
- 2. When the covariances are equal to zero, the symmetric matrix to the 1/2 power is simply the square root of the non-zero entries of the 2 X 2 matrix.

3. In this case  $[V]^{1/2}$  is evaluated by using the matrix formula

$$[V]^{1/2} = [S]^{-1}[\Lambda]^{n}[S]$$
 (22)

where  $[\Lambda]$  and [S] are matrices that diagonalize [V], and the diagonal elements of  $[\Lambda]$  are the eigenvalues of [V] (Ref 3:244-251).

4. To generate samples from the standard normal distribution the central limit theorem was used where the distribution

$$t = \frac{x - \mu}{\frac{\sigma}{\sqrt{n}}}$$
 23)

approaches the standard normal distribution as  $n \to \infty$ . If t = x - 1/2 where x is a pseudo-random number from the uniform distribution, it is found that for a sample of 12 numbers

$$Y = \sum_{i=1}^{12} x_i - 6$$
 (24)

has a distribution that approaches the standard normal distribution where the x<sub>i</sub> are pseudo-random numbers from the uniform distribution (Ref 9:145, 146). Function subprogram RANOR accomplishes the above and returns the numbers to SYSRE.

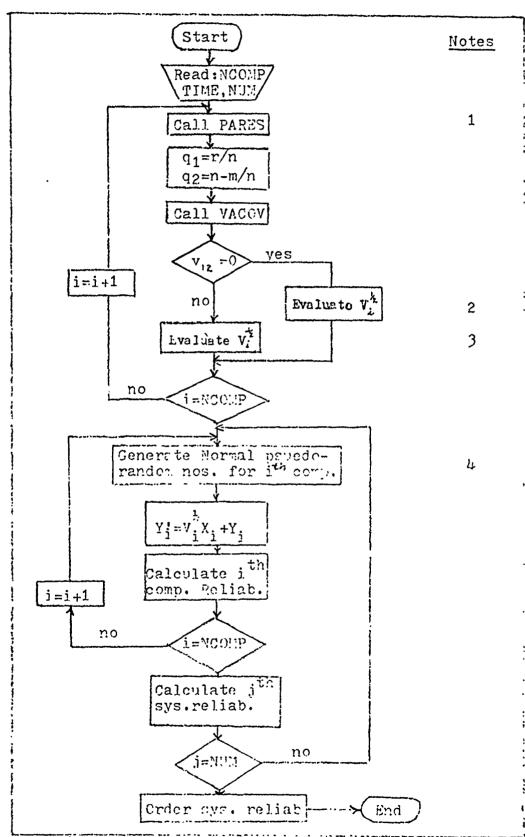


Fig. 4. Flowchart for Main Program

```
PRCGRAM SYSPF(INPUT,OUTPUT)

C*****NCCMP=NUMPER OF CCMCONENTS IN SYSTEM

C*****NUM=NUMPER OF SIMULATIONS DESIRED

C*****TIME=MISSION TIME

C*****FEK=ESIIMATE OF SCALE PARAMETER FOR THAT COMPONENT

C*****ETH=ESTIMATE OF SCALE PARAMETER FOR THAT CCMFONENT

C*****XC=LCCATION FARAMETER FOR THAT CCMFONENT

C*****MABOURGUATION PARAMETER GENSOPING(N=F)C CP LESS)

C*****MABOURGUATION MHEEP OF SAMPLES CENSOPING (N=F)C CP FROM ABOVE

C*****MRLCW=KUMPER OF SAMPLES CENSOPING FROM BFLOW

OTMFNSION S(2,2),V(131),PRLOW(100),EEK(13),FTH(10),

CNSAMP(130),*VA3OV(131),PRLOW(100),EEK(13),FTH(10),

RS(130)=0.

RS(1300)=0.

RS(1300)=0.
                         PROGRAM SYSPF (INPUT, OUTPUT)
            RS(1000)=0.
RS(1000)=0.
RS(1000)=0.
READ 50, NCCYP, NUM, TIME
FO FORMAT(15, 15, F20. P)
       DI = AA/CC
                         US=(CC-65)
                         CÄLL VAČCV(Č1,G2,5K(K),CC,TH(K),V11,V12,V22)
                         A=V11
                         2=V12
                        E=V12

C=V2?

IF (V12)11[,170,117

V(K,1,1)=SGRT(V11)

V(K,1,2)=C.

V(K,2,1)=C.

V(K,2,2)=SCRT(V22)

V(1,2)=C.

V(2,1)=C.

V(2,2)=SCRT(V22)

GO TC 112

Y=.5*CQRT(A*A-2.*A*C+C*C+4.*E*P)

X1=.5*(A+C)+Y
                        Y=.5*SQPT(A*A-2.*A*C+C*C+4

X1=.5*(A+C)+Y

X?=.5*(A+C)+Y

R1=SCPT(1.+(X1-A)**2/B**2)

R2=SQPT(1.+(X2-C)**2/P**2)

S(1,1)=1./F1

S(2,2)=1./F2

S(1,2)=(X2-C)/(P*P2)

S(2,1)=(Y1-A)/(B*P1)

X1=SCPT(Y2)
                        X1=5(0+1(X1)
X2=5(0+1(X2))
D1=5(1,1)*X1
D2=5(1,2)*X2
D3=5(2,1)*X1
C4=5(2,2)*X2
```

```
V(K,1,1)=C1*S(1,1)+C2*S(1,2)
V(K,1,2)=C1*S(2,1)+C2*S(1,2)
V(K,2,1)=C3*S(2,1)+C4*S(1,2)
V(K,2,2)=C3*S(2,1)+C4*S(1,2)
V(K,2,2)=C3*S(2,1)+C4*S(1,2)
V(K,2,2)=C3*S(2,1)+C4*S(1,2)
V(K,2,2)=C3*S(2,1)+C4*S(1,2)
V(K,2,2)=C3*S(2,1)+C4*S(1,2)
V(K,2,2)=C3*S(2,1)+C4*S(1,2)
V(K,2,2)=C3*S(2,1)+C4*S(1,2)
V(K,2,2)
I13 POINT 11E
V(K,2,1)+C4*S(1,2)
I13 POINT 11E
V(K,1,1)+C4*S(1,2)
V(K,2,2)
V(K,2,1)+C4*S(1,2)
V(K,2,2)
V(K,2,
                                                                                                                                           C******CONTINUE

C*****GENERATING SAMPLE SYSTEM RELIGILITY

C*****ORCERING SAMPLE SYSTEM PELIABILITY

ML=1

RING SAMPLE SYSTEM PELIABILITIES
                                                                                                                                                                                            MHT=WH

ISW=1

DO 200

IF ($\overline{\text{T}}$ \text{T} \text{T}
                                    1er Isk=2

                                                                                                                                                                                                                                                                                                                                  , NIN , *CYSTEM RELIABILITY SAMPLE, N=*, 14,
                        J3=J+23n
270 J4=2+4Jn
CPS(J4) ?75,J,R9(J),J1,PS(J1),J2,RS(J2),J3,R5(J3),J4,
270 PPINT 275, J, R9 (J), J1, P9 (J1),
275 PORMAT (1H, 5: [14, 2X, F6.4, 2Y))
300 IF (NUM. Nc. cap) GC TO 40n
J2=J+2:0;
J2=J+40;
J3=J+60;
J4=J+800
```

```
213 FRIRT 315, J,RS(J), J1,RS(J1), J2,RS(J2), J3,RS(J3), J4,

315 FORMAT(1H, 5(I4,2x,F5.4))

400 IF(NJM.NE. 2009)RD TC 5CC

J1=J+ED

J2=J+10JD

J4=J+20JD

J4=J+20JD

J4=J+20JD

J5=J+10JD

J8=J+10JD

FUNCTION RAMOR(Y)

J=D

SUM=CONTINUE

FUNCTION RAMOR(Y)

J=D

SUM=CONTINUE

SUM=CONTINUE

Y=5UM=CONTINUE

PARCH (Y)

TO CONTINUE

PARCH (Y)

PARCH (Y)

PARCH (Y)

PARCH (Y)

PREJUNA

END
```

### Appendix C

# The Variance-Covariance Matrix

Subroutine VACOV solves Eqs (13), (14), and (15); inverts the result and multiplies these by the appropriate constants as shown in Eq (17). The function subprograms GAMI, DGAMI, and D2GAMI compute values of the incomplete Gamma function and its derivatives by using a series expansion as shown in Ref 6. The function subprograms GAM, DGAM and D2GAM use the asymptotic expansion formula to calculate values of the complete Gamma function and its derivatives (Ref 4:257). The flowchart for subroutine VACOV is given in Fig. 5.

# Notes (see Fig. 5):

1. 
$$v_{11} = -Kp + K(K+1)[\Gamma(2;\hat{z}_m^K)] + K(K+1)q_2\hat{z}_m^K$$
 (25)

$$v_{22} = \frac{p}{\kappa^2} + \frac{[\Gamma''(2; \hat{z}_m^K)]}{\kappa^2} + q_2 \hat{z}_m^K \ln^2 \hat{z}_m$$
 (26)

$$v_{12} = p - [\Gamma'(2; \hat{z}_m^K)] - [\Gamma(2; \hat{z}_m^K)] -$$

$$q_2 \hat{z}_m^K [1 + K \ln \hat{z}_m^K]$$
 (27)

2. 
$$v_{11} = -Kp + K(K+1)$$
 (28)

$$v_{22} = \frac{p}{\kappa^2} + \frac{\Gamma''(2.0)}{\kappa^2}$$
 (29)

$$v_{11} = v_{11} + [-K(K+1)]\Gamma(2.0; \hat{z}_{r+1}^{K}) + \frac{q_{1}\hat{z}_{r+1}f(\hat{z}_{r+1})[K\hat{z}_{r+1}^{K} - (K+1)q_{1}]}{q_{1}^{2}}$$

$$v_{22} = v_{22} - \frac{\Gamma''(2.0; \hat{z}_{r+1}^{K})}{\sqrt{2}} + \frac{q_{1}\hat{z}_{r+1}f(\hat{z}_{r+1})[\hat{z}_{r+1}^{K} - q_{1}]}{q_{1}^{2}K}$$

$$v_{12} = v_{12} + \Gamma'(2.0; \hat{z}_{r+1}^{K}) + \Gamma(2.0; \hat{z}_{r+1}^{K}) - \frac{q_{1}\hat{z}_{r+1}f(\hat{z}_{r+1})[K\hat{z}_{r+1}^{K} + q_{1}]}{q_{1}^{2}K}$$

$$v_{13} = v_{14} + \Gamma'(2.0; \hat{z}_{r+1}^{K}) + \Gamma(2.0; \hat{z}_{r+1}^{K}) - \frac{q_{1}\hat{z}_{r+1}f(\hat{z}_{r+1})[K\hat{z}_{r+1}^{K} + q_{1}\hat{z}_{r+1} - (K \ln \hat{z}_{r+1} + 1)q_{1}]}{q_{1}^{2}K}$$

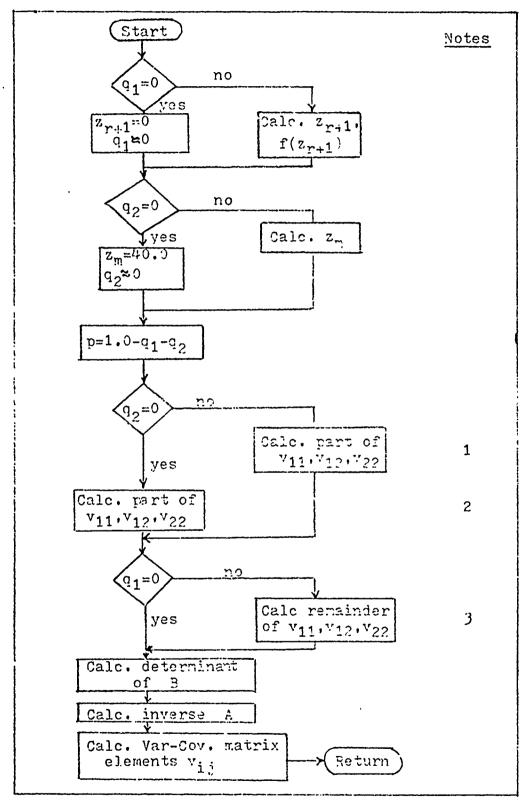


Fig. 5. Flowchart for the Variance-Covariance Subroutine

```
SUBPOUTING MACCV (31,02,07,5MUM,TH,511,512,522)
 X=3
J=3
TF(31.20.1.)GC TC 1
TF(62.50.1.)GC TC 1
GO TO 3
1 PRINT 3
2 FORMAT(14 ,144ALL !! CENSORED)
GO TO 27
    ͱή
                                                   Best odiced from SOPY.
 3 CONTINUT

IF (01.20...)60 TO (

7k=1-1106(1.00-11)1**(1.../2K)

FLZ9=5K*76**(:K-1...)* TXD(-76**5K)

60 TO 5

4 ZR= .6

V=1

01-1 -22
 01=1.L-23
5 IF(02.F0.1.160 TO 7
74=1-HL03(02))**(1.,7/54)
60 TO 3
   74=4 ...
```

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Seprodiced from Solis of Solis
                                                                                         FUNCTION EGAM (Y)

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126-1-1.1.2.7

127-1-1.1.2.7

128-1-1.1.2.7

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     3
                                                                                                                             FUNCTION GAMI(W,7)
DIVENSTON ((F))
U(1)=W**Z/?
FUEL(1)
DI 1 LE2,57
LLTL-1
                                                                                                                     LLTL-1
ELLTEL
U(L)=(+U(LL)/TLL)*U*(7+FLL-1._.)/(7+FLL)
GL*(=CU
ELTE-1
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FUNCTION

OIM ACTON

U(1) = 1/4 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/2 × 1/
```

### Appendix D

# Generating Sample Component Times to Failure

The program in this appendix was used to generate sample failure times for each component. If the points in the cumulative distribution function are represented by random numbers in the uniform distribution over the interval 0.0 to 1.0 (represented by Y) then

$$Y = F(t) = 1 - exp \left[ -\left(\frac{t-C}{\theta}\right)^{K} \right]$$
 (34)

for the Weibull distribution. The inverse relation is

$$t = -\theta[-\ln(1-Y)]^{1/K} + C$$
 (35)

where t is a random value from the Weibull distribution and Y is a pseudo-random number from the above uniform distribution (Ref 15:258). Program RSAMP calculates the true component and system reliability for the given number of components using the required mission time. It then generates the individual component failure times and orders them so that the order statistics can be read into the subroutine which calculates the MLE for shape and scale parameters. This program was used to check the accuracy of the computed confidence interval and confidence level and is not a part of the method proposed.

```
12 FUNCTION TO K (1/2 ) TO COMPONENTS = ,12, 15 HM ISSICH TIME = ,

12 FUNCTION TO K (2/2) TO COMPONENTS = ,12, 15 HM ISSICH TIME = ,
                              12 FURNAT(11) , 21HNUMBER OF COMPONENTS=,12,12HNISSICN TIME=,

CF8.4,7/)

DC. ?) K=1, knowe

15 FORMAT(7HLL, G, 3TT)

CONTINUE

CONTINUE
                                 SUBTRUTING PARET (TV, TH, YC, T, N, K)

2 FORMAT (14 .59 (*.*), //, 17 , 17 HOOMPONINT YUMBER , IN, //)

4 FORMAT (14 , X, EMTHTIA, 7X, 1HX, 3X, 4HLCC., /, 1H , 3ER. 4,
                                              Z=-(ALOO(1.^-Y))
T(T)=(TF*7**(1. /=Y))+;
                 CONTINUE
                                                                                                                                                                                                                                                                                                       Reproduced from sociable con
                                  MPT=N-1

TCW=1

DO CT I=ML, WHT

IF (T(I)-T(T+1)) CL, SE, 2E

GO TO (SE, 45), ISW

TEMPET (T)

T(I)=T(T+1)

T(T+1)=TEME
                             T(T+1)=TEVE
CONTINUE
GC TC(55, 20), TSV
CONTINUE
CONTINUE
```

ŧ

### Vita

Robert Grantham Lannon, the son of Robert E. and Marian G. Lannon, was born on 28 October 1940, in Syracuse, New York. He was graduated from Jamesville-DeWitt High School, Lyndon, New York, in 1958 and entered the United States Air Force Academy that same year. He was graduated from the academy in June 1962 with a bachelors degree in Military Science and a commission as a 2nd Lieutenant in the USAF. After attending pilot training, he served in a variety of flying assignments before entering the Air Force Institute of Technology in June 1970.

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